



Form PTO-1449 Modified List of Patent and Publications Cited by Applicant (Use several sheets if necessary) U.S. Department of Commerce Patent and Trademark Office		Docket No. IBIS-0339	Serial No. 09/753,869
		Applicant Richard Griffey and Eric Swayze	
		Filing Date January 3, 2001	Group Unknown
OTHER DOCUMENTS (Including Author, Title, Date, Pertinent Pages, Etc.)			
<i>MB</i>	AA	Atherton, et al., in <i>Solid Peptide Synthesis: A practical Approach</i> , IRL Press, Oxford, UK, 1989 , 135	
	AB	Brennan, T., et al., "Two-dimensional parallel array technology as a new approach to automated combinatorial solid-phase organic synthesis," <i>Automation</i> , 1998 , 33-45	
	AC	Chen et al., "Structure-Based Discovery of Ligands Targeted to the RNA Double Helix", <i>Biochem.</i> , 1997 , Vol. 36, pp. 11402-11407	
	AD	Chen et al., "Spectroscopic Recognition of Guanine Dimeric Haripin Quadruplexes by a Carbocyanine Dye", <i>Proc. Natl. Acad. Sci.</i> , 1996 , Vol. 93, pp. 2635-2639	
	AE	Corey, E.J. et al., "Computer-Assisted Analysis in Organic Synthesis," <i>Science</i> , 1985 , 228, 408-418	
	AF	Czarnik, A.W. et al. (eds.), <i>A Practical Guide to Combinatorial Chemistry</i> , American Chemical Society, Washington, DC, 1997 , Chs. 2, 13, and 14, 17-47 and 357-412	
	AG	Gschwend et al., "Orientational Sampling and Rigid-Body Minimization in Molecular Docking Revisited: On-the-fly Optimization and Degeneracy Removal", <i>J. Comput.-Aided Mol. Des.</i> , 1996 , Vol. 10, pp. 123-132	
	AH	Heathcock, et al., "Total synthesis of racemic vallesamidine," <i>J. Org. Chem.</i> ®, 1990 , 55(3), 798-911 (abstract only)	
	AI	Kuntz et al., "A Geometric Approach to Macromolecule-Ligand Interactions", <i>J. Mol. Biol.</i> , 1982 , Vol. 161, pp. 269-288	
	AJ	Kuntz et al., "Structure-Based Molecular Design", <i>Acc. Chem. Res.</i> , 1994 , Vol. 27, No. 5, pp. 117-123	
	AK	Levine et al., "Stalk: An Interactive System for Virtual Molecular Docking", <i>IEEE Computational Science and Engineering</i> , pp. 55-65	
	AL	Meng et al., "Automated Docking with Grid-Based Energy Evaluation", <i>J. Comput. Chem.</i> , 1992 , Vol. 13, No. 4, pp. 505-524	
<i>MB</i>	AM	Rotstein, S.H. et al., "GroupBuild: A Fragment-Based Method for <i>De Novo</i> Drug Design," <i>J. Med. Chem.</i> , 1993 , 36, 1700-1710	
EXAMINER		DATE CONSIDERED	



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U. S. PATENT DOCUMENTS

Examiner Initial		Document No.	Date	Name	Class	Subclass
<i>MM</i>	AP	5,010,175	04/23/91	Rutter, et al.	530	334
	AQ	5,186,898	02/16/93	Bridgham, et al.	422	102
	AR	5,372,672	12/13/94	Seifert, et al.	156	584
	AS	5,424,186	06/13/95	Fodor, et al.	435	006
	AT	5,434,796	07/18/95	Weininger	364	496
	AU	5,472,672	12/05/95	Brennan	422	131
	AV	5,529,756	06/25/96	Brennan	422	131
	AW	5,565,324	10/1996	Still et al.	435	6
	AX	5,573,905	11/1996	Lerner et al.	435	6
	AY	5,574,656	11/12/96	Agrafiotis et al.	364	500
	AZ	5,703,792	12/30/97	Chapman	364	496
<i>MM</i>	BA	5,880,972	03/09/99	Horibeck	364	496

FOREIGN PATENT DOCUMENTS

Examiner Initial		Document No.	Date	Country	Translation YES NO	

EXAMINER <i>MM</i>	DATE CONSIDERED <i>02/02</i>
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